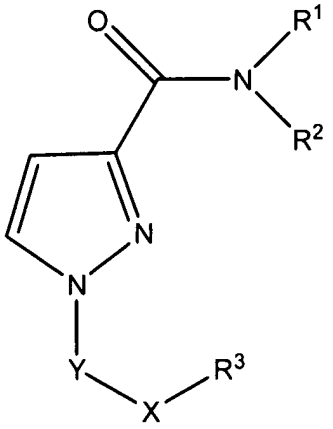
Amendments to the Claims

Please amend Claims 1, 16, 18, 37, 40, 42, 48 and 53. The Claim Listing below will replace all prior versions of the claims in the application:

Claim Listing

1. (Currently Amended) A compound of formula I,



I

wherein either R^1 represents an aryl group or a heteroaryl group, both of which are optionally substituted by one or more substituents selected from G^1 and B^1, which B^1 group may itself be further substituted by one or more substituents selected from G^2, Z (provided that Z is not directly attached to an aryl or a heteroaryl group) and B^2 (which B^2 group is optionally further substituted by one or more substituents selected from G^3, B^3 and Z, provided that Z is not attached to an aryl or a heteroaryl group); and

R^2 represents H or C_{1-6} alkyl, which latter group is optionally substituted by one or more halo groups;

or

when R^2 represents C_{1-6} alkyl optionally substituted by halo, R^1 and R^2 may be linked together forming a further 5- to 7-membered ring, optionally containing 1 to 3 heteroatoms and/or 1 to 3 double bonds, which ring is itself optionally substituted by one

or more substituents selected from G^1, Z (provided that the ring is not aromatic in nature) and B^1 (which B^1 group is optionally substituted as described above);

R^3 represents C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{3-8} cycloalkyl, C_{3-8} heterocycloalkyl, aryl or heteroaryl, all of which groups are optionally substituted by one or more substituents selected from G^{1a}, Z (provided that Z is not directly attached to an aryl or a heteroaryl group) and B^1 (which B^1 group is optionally substituted as described above);

X represents a direct bond, -O- or -N(R^4)-;

Y represents -C(O)-, -C(S)- or -$S(O)_2$-;

B^1, B^2 and B^3 independently represent, on each occasion when used above, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{3-8} cycloalkyl, C_{3-8} heterocycloalkyl, aryl or heteroaryl;

G^1, G^{1a}, G^2 and G^3 independently represent, on each occasion when used above, halo, cyano, -N_3, -NO_2, -ONO_2 or -A^1 -R^4;

wherein A^1 represents a spacer group selected from -C(Z)A^2-, -N(R^5)A^3-, -OA^4-, -S- or -$S(O)_n$$A^5$-, in which:

A^2 represents a single bond, -O-, -S- or -N(R^5)-;

A^3 represents A^6, -C(Z)N(R^5)C(Z)N(R^5)-, -C(Z)N(R^5)C(Z)O-, -C(Z)N(R^5)$S(O)_n$N(R^5), -C(Z)S-, -$S(O)_n$–, -$S(O)_n$N(R^5)C(Z)N(R^5)-, -$S(O)_n$N(R^5)C(Z)O-, -$S(O)_n$N(R^5)$S(O)_n$N(R^5)-, -C(Z)O-, -$S(O)_n$N(R^5)- or -$S(O)_n$O-;

A^4 represents A^6, -$S(O)_n$-, -C(Z)O-, -$S(O)_n$N(R^5)- or -$S(O)_n$O-;

A^5 represents a single bond, -N(R^5)- or -O-;

A^6 represents a single bond, -C(Z)- or -C(Z)N(R^5)-;

Z represents, on each occasion when used above, a substituent connected by a double bond, which is selected from =O, =S, =NR^4, =NN(R^4)(R^5), =NOR^4, =$NS(O)_2$N(R^4)(R^5), =NCN, =$CHNO_2$ and =C(R^4)(R^5);

R^4 and R^5 independently represent, on each occasion when used above, H or B^4, which B^4 group is itself optionally substituted by one or more substituents selected from G^4, Q provided that Q is not directly attached to an aryl or a heteroaryl group) and B^5 (which B^5 group is itself optionally substituted by one or more substituents selected from G5, Q (provided that Q is not directly attached to an aryl or a heteroaryl group) and B^6; or

when R^4 and R^5 both represent optionally substituted B^4 groups, then any pair thereof may, for example when present on the same atom or on adjacent atoms, be linked together to form, with those, or other relevant, atoms, a 5- to 7-membered ring, optionally containing 1 to 3 heteroatoms and/or 1 to 3 double bonds, which ring is itself optionally substituted by one or more substituents selected from G^6, Q (provided that the ring is not aromatic in nature) and B^4 (which B^4 group is optionally substituted as described above);

B^4, B^5 and B^6 independently represent on each occasion when used above C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{3-8} cycloalkyl, C_{3-8} heterocycloalkyl, aryl or heteroaryl;

G^4, G^5 and G^6 independently represent on each occasion when used above, halo, cyano, N_3, $-NO_2$, $-ONO_2$ or $-A^7-R^6$;

wherein A^7 represents a spacer group selected from $-C(Q)A^8-$, $-N(R^7)A^9-$, $-N(R^{7a})A^{9a}-$, $-OA^{10}-$, -S- or $-S(O)_nA^{11}-$, in which:

A^8 represents a single bond, -O-, -S- or $-N(R^7)-$;

A^9 represents A^{12}, $-C(Q)S-$, $-S(O)_n-$, $-C(Q)O-$, $-S(O)_nN(R^7)-$ or $-S(O)_nO-$;

A^{9a} represents $-C(Q)N(R^7)C(Q)N(R^7)-$, $-C(Q)N(R^7)C(Q)O-$, $-C(Q)N(R^7)S(O)_nN(R^7)-$, $-S(O)_nN(R^7)C(Q)N(R^7)-$, $-S(O)_nN(R^7)C(Q)O-$, $-S(O)_nN(R^7)S(O)_nN(R^7)-$;

A^{10} represents A^{12}, $-S(O)_n-$, $-C(Q)O-$, $-S(O)_nN(R^7)-$ or $S(O)_nO-$;

A^{11} represents a single bond, $-N(R^7)-$ or -O-;

A^{12} represents a single bond, $-C(Q)-$ or $-C(Q)N(R^7)-$;

Q represents, on each occasion when used above, a substituent connected by a double bond, which is selected from =O, =S, $=NR^6$, $=NN(R^6)(R^7)$, $=NOR^6$, $=NS(O)_2N(R^6)(R^7)$, =NCN, $=CHNO_2$ and $=C(R^6)(R^7)$;

R^6, R^7 and R^{7a} independently represent, on each occasion when used above, H, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{3-8} cycloalkyl, C_{3-8} heterocycloalkyl, aryl or heteroaryl, which latter seven groups are optionally substituted by one or more groups selected from halo, C_{1-6} alkyl (optionally substituted by one or more halo groups) - $N(R^8)R^9$, $-OR^8$, $-ONO_2$ and $-SR^8$; or

provided that they do not represent H, any pair of R^6 and R^7 may, for example when present on the same atom or on adjacent atoms, be linked together to form, with

those, or other relevant, atoms, a 5- to 7-membered ring, optionally containing 1 to 3 heteroatoms and/or 1 to 3 double bonds, which ring is itself optionally substituted by one or more groups selected from halo, C_{1-6} alkyl (optionally substituted by one or more halo groups), $-N(R^8)R^9$, $-OR^8$, $-ONO_2$ and $-SR^8$;

R^8 and R^9 independently represent, on each occasion when used above, H or C_{1-6} alkyl, which latter group is optionally substituted by one or more halo groups; and n represents, on each occasion when used above, 1 or 2;

or a pharmaceutically-acceptable salt thereof,

provided that, when R^2 represents H, Y represents -C(O)- and:

(A) X represents a direct bond and:

i) R^3 represents phenyl, then R^1 does not represent phenyl, 2-methoxyphenyl, 2-thiazolyl or 6-methyl-2-pyridinyl;

ii) R^3 represents 4-fluorophenyl, then R^1 does not represent 2-carbomethoxyphenyl, 3-carbomethoxyphenyl or 2,4-dimethylphenyl;

iii) R^3 represents 2-chlorophenyl, then R^1 does not represent phenyl, 3-bromophenyl or 4-bromophenyl;

iv) R^3 represents 3-chlorophenyl, then R^1 does not represent phenyl, 2-fluorophenyl, 2-chlorophenyl, 2,3-dichlorophenyl or 2,5-dichlorophenyl;

v) R^3 represents 4-chlorophenyl, then R^1 does not represent 3-bromophenyl or 4-methoxyphenyl;

vi) R^3 represents 3-iodophenyl, then R^1 does not represent 2-methoxyphenyl or 2,4-dimethylphenyl;

vii) R^3 represents ~~2,4-dichlororphenyl~~ 2,4-dichlorophenyl, then R^1 does not represent 4-chlorophenyl or 2,3-dichlorophenyl;

viii) R^3 represents 3,5-dinitrophenyl, then R^1 does not represent 2,3-dichlorophenyl;

ix) R^3 represents 2,4-dimethyl-6-oxo-6*H*-pyran-3-yl, then R^1 does not represent 3-carbomethoxyphenyl;

x) R^3 represents methyl, then R^1 does not represent 3,4-dichlorophenyl, 2-methoxyphenyl, 2-thiazolyl, 4-methyl-2-pyridinyl, 6-methyl-2-pyridinyl or 4-acetylphenyl;

xi) R^3 represents ethyl, then R^1 does not ~~reporesent~~ represent phenyl, 2,3-dichlorophenyl, 4-methoxyphenyl, 2-carbomethoxy-phenyl, 2-thiazolyl or 4-methyl-2-pyridinyl;

(B) X represents -N(H)- and:

i) R^3 represents phenyl, then R^1 does not represent 4-methoxyphenyl, 2,4-dimethylphenyl or 2-thiazolyl;

ii) R^3 represents 3-chlorophenyl, then R^1 does not represent 4-methylphenyl;

iii) R^3 represents 4-chlorophenyl, then R^1 does not represent 3-bromophenyl;

iv) R^3 represents 3,4-dichlorophenyl, then R^1 does not represent 4-methyl-2-pyridinyl or 6-methyl-2-pyridinyl;

v) R^3 represents 2'-sulfamoylbiphenyl-4-yl, then R^1 does not represent 5-bromo-2-pyridinyl;

vi) R^3 represents 1-propyl, then R^1 does not represent phenyl;

vii) R^3 represents 1-butyl, then R^1 does not represent 4-bromophenyl or 2,4-dimethylphenyl;

viii) R^3 represents cyclohexyl, then R^1 does not represent 4-methoxyphenyl;

(C) X represents -O- and:

i) R^3 represents phenyl, then R^1 does not represent phenyl or 6-methyl-2-pyridinyl;

ii) R^3 represents methyl, then R^1 does not represent phenyl, 2-fluorophenyl, 2,4-dimethylphenyl, 4-acetylphenyl or 2-thiazolyl;

iii) R^3 represents ethyl, then R^1 does not represent phenyl, 2-fluorophenyl, 4-acetylphenyl or 4-methyl-2-pyridinyl;

iv) R^3 represents 1-butyl, then R^1 does not represent 2-fluorophenyl, 2-methoxyphenyl, 4-methyl-2-pyridinyl or 6-methyl-2-pyridinyl;

v) R^3 represents 2-butyl, then R^1 does not represent 2-thiazolyl or 4-acetylphenyl;

vi) R^3 represents 2-methyl-1-propyl, then R^1 does not represent phenyl or ~~3-nitorphenyl~~ 3-nitrophenyl.

2. (Original) A compound as claimed in Claim 1, wherein R^1 represents an aryl or heteroaryl group, both of which are optionally substituted as defined in Claim 1.

3. (Previously Presented) A compound as claimed in Claim 1, wherein G^1 represents halo, cyano or $-A^1-R^4$.

4. (Previously Presented) A compound as claimed in Claim 1, wherein B^1 represents an optionally substituted C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{3-7} cycloalkyl, C_{4-7} heterocycloalkyl, or phenyl, group.

5. (Previously Presented) A compound as claimed in Claim 1, wherein G^{1a} represents halo, cyano, $-NO_2$ or $-A^1-R^4$.

6. (Previously Presented) A compound as claimed in Claim 1, wherein G^2 represents halo, cyano, $-ONO_2$ or $-A^1-R^4$.

7. (Previously Presented) A compound as claimed in Claim 1, wherein B^2 represents C_{1-6} alkyl, C_{2-6} alkenyl or C_{2-6} alkynyl, all of which are optionally substituted by one or more G^3 and/or B^3 groups.

8. (Previously Presented) A compound as claimed in Claim 1, wherein G^3 represents halo, $-ONO_2$, $-N(R^5)(R^4)$ or $-OR^4$.

9. (Previously Presented) A compound as claimed in Claim 1, wherein B^3 represents C_{1-6} alkyl, C_{2-6} alkenyl or C_{2-6} alkynyl.

10. (Previously Presented) A compound as claimed in Claim 1, wherein when A^1 represents $-N(R^5)A^3-$, A^3 represents A^6, $-C(Z)S-$, $-S(O)_n-$, $-C(Z)O-$ or $-S(O)_nN(R^5)-$.

11. (Previously Presented) A compound as claimed in Claim 1, wherein when A^1 represents $-OA^4-$, A^4 represents A^6.

12. (Previously Presented) A compound as claimed in Claim 1, wherein when A^1 represents $-S(O)_nA^5-$, A^5 represents a single bond or $-N(R^5)-$.

13. (Previously Presented) A compound as claimed in Claim 1, wherein when A^1 represents $-C(Z)A^2-$, A^2 represents a single bond, -O- or $-N(R^5)-$.

14. (Previously Presented) A compound as claimed in Claim 1 wherein A^1 represents $-C(Z)A^2-$, $-N(R^5)A^3-$ or $-OA^4-$.

15. (Previously Presented) A compound as claimed in Claim 1, wherein Z represents =O or $=NR^4$.

16. (Currently Amended) A compound as claimed in Claim 1, wherein when any pair [[or]] <u>of</u> R^4 and R^5 are linked together to form a ring, they are optionally substituted with G^6 and/or B^4.

17. (Previously Presented) A compound as claimed in Claim 1, wherein G^4 represents halo, cyano, $-ONO_2$ or $-A^7-R^6$.

18. (Currently Amended) A compound as claimed in Claim 1, wherein B^5 represents C_{1-6} alkyl, C_{2-6} alkenyl or C_{2-6} <u>alkynyl</u> ~~alknyl~~, all of which are optionally substituted by one or more G^5 and/or B^6 groups.

19. (Previously Presented) A compound as claimed in Claim 1, wherein G^5 represents halo, $-ONO_2$, $-N(R^7)(R^6)$ or $-OR^6$.

20. (Previously Presented) A compound as claimed in Claim 1, wherein B^6 represents C_{1-6} alkyl, C_{2-6} alkenyl or C_{2-6} alkynyl.

21. (Previously Presented) A compound as claimed in Claim 1, wherein G^6 represents halo, cyano or $-A^7-R^6$.

22. (Previously Presented) A compound as claimed in Claim 1, wherein A^7 represents $-C(Q)A^8-$, $-N(R^7)A^9-$, $-OA^{10}-$, -S- or $-S(O)_nA^{11}-$.

23. (Previously Presented) A compound as claimed in Claim 1, wherein when A^7 represents $-N(R^7)A^9-$, A^9 represents A^{12}, $-C(Q)S-$, $-S(O)_n-$, $-C(Q)O-$ or $-S(O)_nN(R^7)-$.

24. (Previously Presented) A compound as claimed in Claim 1, wherein when A^7 represents $-OA^{10}-$, A^{10} represents A^{12}.

25. (Previously Presented) A compound as claimed in Claim 1, wherein when A^7 represents $-S(O)_nA^{11}-$, A^{11} represents a single bond or $-N(R^7)-$.

26. (Previously Presented) A compound as claimed in Claim 1, wherein when A^7 represents $-C(Q)A^8-$, A^8 represents a single bond, -O- or $-N(R^7)-$.

27. (Previously Presented) A compound as claimed in Claim 1, wherein Q represents =O or $=NR^6$.

28. (Previously Presented) A compound as claimed in Claim 1, wherein R^6, R^7 and R^{7a} independently represent H, C_{1-6} alkyl, C_{2-6} alkenyl or C_{2-6} alkynyl, all of which groups are optionally substituted by one or more groups selected from halo, C_{1-6} alkyl, $-N(R^8)R^9$, OR^8 and $-ONO_2$.

29. (Previously Presented) A compound as claimed in Claim 1 wherein when any pair of R^6 and R^7 are linked together to form a ring, that ring is optionally substituted by one or more groups selected from halo, C_{1-6} alkyl (optionally substituted by one or more halo groups), $-N(R^8)R^9$, $-OR^8$ and $-ONO_2$.

30. (Previously Presented) A compound as claimed in Claim 1, wherein B^4 represents an optionally substituted C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{3-7} cycloalkyl, C_{4-7} heterocycloalkyl, or phenyl, group.

31. (Previously Presented) A compound as claimed in Claim 1 wherein R^4 and/or R^5 independently represent H or C_{1-6} alkyl, which latter group is optionally substituted by one or more fluoro groups.

32. (Previously Presented) A compound as claimed in Claim 1, wherein X represents a direct bond, -O-, -N(H)- or -N(Me)-.

33. (Previously Presented) A compound as claimed in Claim 1 wherein R^2 represents H, methyl or ethyl.

34. (Previously Presented) A compound as claimed in Claim 1, wherein R^1 represents an optionally substituted phenyl, naphthyl, pyrrolidinyl, piperidinyl, pyrrolyl, furanyl, thiophenyl, pyrazolyl, imidazolyl, oxazolyl, isoxazolyl, thiazolyl, pyridinyl, indazolyl, indolyl, indolinyl, isoindolinyl, oxindolyl, quinolinyl, 1,2,3,4-tetrahydroquinolinyl, isoquinolinyl, 1,2,3,4-tetrahydroisoquinolinyl, quinolizinyl, benzofuranyl, isobenzofuranyl, chromanyl, benzothiophenyl, pyridazinyl, pyrimidinyl, pyrazinyl, indazolyl, benzimidazolyl, quinazolinyl, quinoxalinyl, 1,3-benzodioxolyl, benzothiazolyl, or benzodioxanyl, group.

35. (Original) A compound as claimed in Claim 34, wherein R^1 represents optionally substituted phenyl, 2-pyridinyl, 3-pyridinyl, 2-thiophenyl, 4-pyrazolyl, 5-isoxazolyl, 1,3-benzodioxolyl, indazolyl, benzothiazolyl, or quinolinyl, group.

36. (Previously Presented) A compound as claimed in Claim 34, wherein the optional substituent(s) are selected from halo, cyano, C_{1-6} alkyl (which alkyl group may be linear or branched, and/or substituted by one or more fluoro and/or C_{3-6} cycloalkyl groups), C_{2-6} alkenyl, C_{3-6} cycloalkyl, phenyl, pyrrolidinyl, piperidinyl, piperazinyl, tetrahydrofuranyl,

tetrahydropyranyl, morpholinyl, thiomethyl, methylsulfinyl, methylsulfonyl, -OR^{10}, -$N(R^{10})R^{11}$, -$C(O)OR^{10}$, -$C(O)R^{10}$, -$C(O)N(R^{10})R^{11}$, -$S(O)_2N(R^{10})R^{11}$ and -$N(R^{10})S(O)_2R^{12}$, wherein R^{10} and R^{11} independently represent H, phenyl, C_{1-6} alkyl (which alkyl group is optionally substituted by one or more fluoro atom), C_{2-6} alkenyl or C_{3-6} cycloalkyl; or R^{10} and R^{11} may be linked together to form, with the nitrogen atom to which they are attached, a 5- to 7-membered ring, optionally containing one additional heteroatom and optionally substituted with one or more C_{1-6} alkyl groups, which alkyl groups are themselves optionally substituted by one or more halo groups; and R^{12} represents phenyl, C_{1-6} alkyl (which alkyl group is optionally substituted by one or more fluoro atom), C_{2-6} alkenyl or C_{3-6} cycloalkyl.

37. (Currently Amended) A compound as claimed in Claim 36, wherein the optional substituent(s) are selected from carbomethoxy, methyl, dimethylamino, cyano, chloro ~~cholor~~, fluoro, trifluoromethyl, bromo, methoxy and trifluoromethoxy.

38. (Previously Presented) A compound as claimed in Claim 1, wherein R^3 represents an optionally substituted C_{1-6} alkyl, C_{3-6} cycloalkyl, phenyl, naphthyl, pyrrolidinyl, piperidinyl, piperazinyl, pyrrolyl, furanyl, thiophenyl, pyrazolyl, imidazolyl, oxazolyl, isoxazolyl, thiazolyl, pyridinyl, indazolyl, indolyl, indolinyl, isoindolinyl, oxindolyl, quinolinyl, 1,2,3,4-tetrahydroquinolinyl, isoquinolinyl, 1,2,3,4-tetrahydroisoquinolinyl, quinolizinyl, benzofuranyl, isobenzofuranyl, chromanyl, benzothiophenyl, pyridazinyl, pyrimidinyl, pyrazinyl, indazolyl, benzimidazolyl, quinazolinyl, quinoxalinyl, 1,3-benzodioxolyl, benzothiazolyl, or benzodioxanyl, group.

39. (Original) A compound as claimed in Claim 38, wherein R^3 represents an optionally substituted C_{1-6} alkyl, cyclohexyl, phenyl, 2-thiophenyl, 2-furanyl, 3-furanyl, 2-pyrrolyl, 1-naphthyl, 4-piperazinyl, 4-piperidinyl, benzofuranyl, or 1,3-benzodioxolyl, group.

40. (Currently Amended) A compound as claimed in Claim 38, wherein the optional substituent(s) are selected from halo, -NO_2, cyano, C_{1-6} alkyl (which alkyl group may be linear or branched, and/or optionally substituted with one or more halo, C_{1-6} alkyl, C_{2-6}

alkenyl and/or C_{3-6} cycloalkyl, groups, which latter three groups are themselves optionally substituted with one or more halo and/or C_{1-6} alkyl groups), C_{2-6} alkenyl (optionally substituted with one or more C_{1-6} alkyl groups), C_{3-6} cycloalkyl (optionally substituted with one or more halo groups), phenyl (optionally substituted with one or more halo groups), pyrrolidinyl, piperidinyl, piperazinyl, tetrahydrofuranyl, tetrahydropyranyl, morpholinyl, thiomethyl, ~~methylsufinyl~~ methylsulfinyl, methylsulfonyl, =O, $-OR^{13}$, $-N(R^{13})R^{14}$, $-C(O)OR^{13}$, $-C(O)R^{13}$, $-C(O)N(R^{13})R^{14}$, $-S(O)_2N(R^{13})R^{14}$ and $-N(R^{13})S(O)_2R^{15}$, wherein R^{13} and R^{14} independently represent H, phenyl, C_{1-6} alkyl (which alkyl group is optionally substituted by one or more fluoro atom), C_{2-6} alkenyl or C_{3-6} cycloalkyl; or R^{13} and R^{14} may be linked together to form, with the nitrogen atom to which they are attached, a 5- to 7- membered ring, optionally containing one additional heteroatom and optionally substituted with one or more C_{1-6} alkyl groups, which alkyl groups are themselves optionally substituted by one or more halo groups; and R^{15} represents phenyl, C_{1-6} alkyl (which alkyl group is optionally substituted by one or more fluoro atom), C_{2-6} alkenyl or C_{3-6} cycloalkyl ~~cycloaklyl~~.

41. (Original) A compound as claimed in Claim 40, wherein the optional substituent(s) are selected from methyl, ethyl, ethoxy, trifluoromethyl, fluoro, chloro, iodo, phenyl, 2-chlorophenyl, 4-chlorophenyl, *n*-pentyl, *i*-propyl, nitro, *t*-butyl, $-CH_2CH=CHC_8H_{17}$, trifluoroacetyl, carbomethoxy, carboethoxy and trifluoromethoxy.

42. (Currently Amended) A compound as claimed in Claim 1, wherein R^1 is phenyl, 2-chlorophenyl, 2-chloro-4-fluorophenyl, 3-chloro-4-fluorophenyl, 2,6-dichlorophenyl, 5-chloro-2-cyanophenyl, 2-fluoro-5-trifluoromethylphenyl, 2-bromo-4-trifluoromethoxyphenyl, 2-methoxy-6-methylphenyl, 3-cyanophenyl, 4-trifluoromethylphenyl, 4-dimethylaminophenyl, 4-carbomethoxyphenyl, 1,3,5-trimethyl-1*H*-pyrazol-4-yl, 3-methylisoxazol-5-yl, 3-pyridinyl, 2-chloro-3-pyridinyl, 3-methyl-2-pyridinyl, 3-carbomethoxythiophen-2-yl or 1,3-benzodioxolyl;

R^2 is hydrogen or methyl;

R^3 is methyl, *n*-butyl, *n*-pentyl, 1-octyl, oleoyl, (1*R*,2*S*,5*R*)-(-)-menthyl, 2-chlorobenzyl, benzyl, phenyl, 3-fluorophenyl, 3-chlorophenyl, 4-chlorophenyl, 2-fluoro-5-iodophenyl, 5-fluoro-2-methylphenyl, ~~4-*tert*-butyl-phenyl~~ 4-*tert*-butylphenyl, 4-pentylphenyl, 3-trifluoromethylphenyl, 4-trifluoromethoxyphenyl, 4-nitrophenyl, ~~2-ethoxyphenyl~~ 2-ethoxyphenyl, 1-naphthyl, 2-furanyl, 2,5-dimethyl-3-furanyl, 2-carbomethoxy-5-furanyl, 1-methyl-1*H*-pyrrol-2-yl, 3-methyl-2-benzofuranyl, 3-methyl-2-thiophenyl, 1(*N*)-methyl-4-piperazinyl, 1(*N*)-(2,2,2-trifluoroacetyl)piperidin-4-yl, ethylhexanoate or 1,3-benzodioxolyl; Y is -C(O)-, -C(S)- or $-S(O)_2-$; and

X is a bond, -N(H)-, -N(Me)-, or -O-.

43. (Cancelled)

44. (Previously Presented) A pharmaceutical formulation including a compound of formula I, as defined in Claim 1, but without the provisos, or a pharmaceutically acceptable salt thereof, in admixture with a pharmaceutically acceptable adjuvant, diluent or carrier.

45. (Cancelled)

46. (Withdrawn) The method as claimed in Claim 49 wherein the lipoxygenase is 15-lipoxygenase.

47. (Withdrawn) The method as claimed in Claim 46, wherein the disease is inflammation and/or has an inflammatory component.

48. (Withdrawn, Currently Amended) The method as claimed in Claim 47 wherein the inflammatory disease is asthma, chronic obstructive pulmonary disease (COPD), pulmonary fibrosis, an allergic disorder, rhinitis, inflammatory bowel disease, an ulcer, inflammatory pain, fever, atherosclerosis, coronary artery disease, vasculitis, pancreatitis, arthritis, osteoarthritis, rheumatoid arthritis, conjunctivitis, iritis, scleritis, uveitis, a wound, dermatitis, eczema, psoriasis, stroke, diabetes, autoimmune diseases, Alzheimer's disease, multiple sclerosis, sarcoidosis, Hodgkin's disease or another malignancy.

49. (Withdrawn) A method of treatment of a disease in which inhibition of the activity of a lipoxygenase is desired and/or required, which method comprises administration of a therapeutically effective amount of a compound of formula I as defined in Claim 1, but without the provisos, or a pharmaceutically-acceptable salt thereof, to a patient suffering from, or susceptible to, such a condition.

50. (Previously Presented) A combination product comprising:

(A) a compound of formula I as defined in Claim 1, but without the provisos; and

(B) another therapeutic agent that is useful in the treatment of inflammation, wherein each of components (A) and (B) is formulated in admixture with a pharmaceutically-acceptable adjuvant, diluent or carrier.

51. (Previously Presented) A combination product as claimed in Claim 50 which comprises a pharmaceutical formulation including the compound of formula I, another therapeutic agent that is useful in the treatment of inflammation, and a pharmaceutically-acceptable adjuvant, diluent or carrier.

52. (Previously Presented) A combination product as claimed in Claim 50 which comprises a kit of parts comprising components:

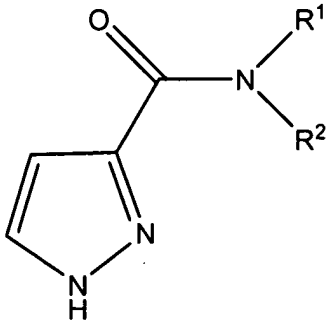
(a) a pharmaceutical formulation including the compound of formula I, in admixture with a pharmaceutically-acceptable adjuvant, diluent or carrier; and

(b) a pharmaceutical formulation including another therapeutic agent that is useful in the treatment of inflammation in admixture with a pharmaceutically-acceptable adjuvant, diluent or carrier,

which components (a) and (b) are each provided in a form that is suitable for administration in conjunction with the other.

53. (Withdrawn, Currently Amended) A process for the preparation of a compound as defined in Claim 1, which comprises:

(i) for compounds of formula I in which, when Y is $-S(O)_2-$, X represents a direct bond or $-N(R^4)-$, in which R^4 represents B^4, reaction of a compound of formula II,



II

wherein R^1 and R^2 are as defined in Claim 1, with a compound of formula III,

$R^3-X^a-Y-L^1$ III

wherein [[X^8]] $\underline{X^a}$ represents a direct bond or $-N(B^4)-$ when Y represents $-S(O)_2-$ or, for all other values of Y, represents X as defined in Claim 1, R^3 and Y are as defined in Claim 1 and L^1 represents a suitable leaving group;

(ii) for compounds of formula I in which X represents a single bond and Y represents -C(O)-, reaction of a compound of formula II as defined above with a compound of formula IV,

$R^3C(O)OH$ IV

wherein R^3 is as defined in Claim 1;

(iii) for compounds of formula I in which X represents a direct bond and Y represents a -C(O)- or a -C(S)- group, reaction of a compound of formula II as defined above with a compound of formula V,

$R^3=Y^a$ V

wherein [[Y^8]] $\underline{Y^a}$ represents -C(O)- or -C(S)- and R^3 is as defined in Claim 1;

(iv) for compounds of formula I, in which X represents -NH- and Y represents -C(O)- or -C(S)-, reaction of a compound of formula II as defined above with a compound of formula VI,

$R^3N=Y^a$ VI

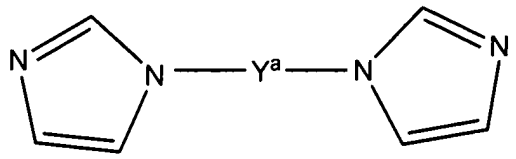
wherein R^3 is as defined in Claim 1 and Y^a is as defined above;

(v) for compounds of formula I in which Y represents -C(O)- or -C(S)-, reaction of a compound of formula II as defined above with:

(a) a compound of formula VII,

~~Cl-Y^a-Cl~~ Cl-Y^a- Cl VII

(b) a compound of formula VIII,



VIII

wherein, in both cases, Y^a[[8]] is as defined above; or

(c) when Y represents -C(O)-, triphosgene, followed by:

(1) for compounds of formula I in which X represents a direct bond, reaction with a compound[[s]] of formula IX,

R^3M IX

wherein M represents a metal such as Mn, Fe, Ni, Cu, Zn, Pd or Ce, or a salt or complex thereof and R^3 is as defined in Claim 1;

(2) for compounds of formula I wherein X represents O, reaction with a compound of formula X,

R^3OH X

wherein R^3 is as defined in Claim 1; or

(3) for compounds of formula I wherein X represents -N(R^4)-, reaction with a compound of formula XI,

$R^3N(H)R^4$ XI

wherein R^3 and R^4 are as defined in Claim 1;

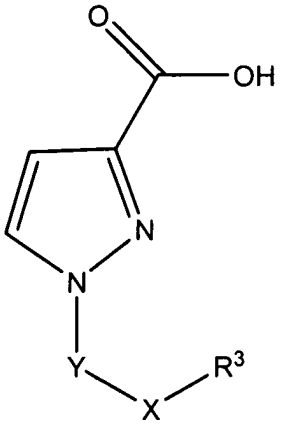
(vi) for compounds of formula I in which X represents -N(R^4)- and R^4 is other than hydrogen, reaction of a corresponding compound of formula I in which X represents -N(H)- with a compound of formula XII,

R^4-L^1 XII

wherein R^4 is as defined in Claim 1 and L^1 is as defined above;

(vii) for compounds of formula I in which Y represents -C(S)-, reaction of a corresponding compound of formula I in which Y represents -C(O)- with a suitable reagent for the conversion of a carbonyl group to a thiocarbonyl group;

(viii) reaction of a compound of formula XIII,



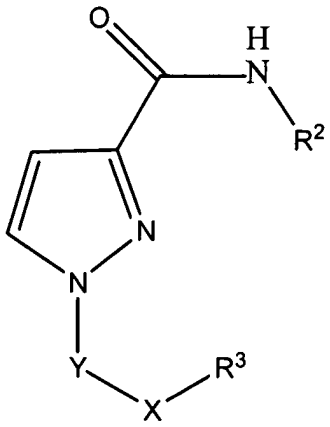
XIII

wherein R^3, Y and X are as defined in Claim 1, with a compound of formula XIV,

$HN(R^1)(R^2)$ XIV

wherein R^1 and R^2 are as defined in Claim 1; or

(ix) reaction of a compound of formula XV,



XV

wherein R^2, R^3, Y and X are as defined in Claim 1, with a compound of formula XVI,

R^1-L^2 XVI

wherein L^2 represents a suitable leaving group and R^1 is as defined in Claim 1.